

Methanimine, 1-(1-pyrrolidiny), N-hexyl

Inchi:	InChI=1S/C11H22N2/c1-2-3-4-5-8-12-11-13-9-6-7-10-13/h11H,2-10H2,1H3/b12-11+
InchiKey:	UPYONVCINBLBAV-VAWYXSNFSA-N
Formula:	C11H22N2
SMILES:	CCCCCCN=CN1CCCC1
Mol. weight [g/mol]:	182.31

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.55		Crippen Method
logp	2.691		Crippen Method
mcvol	170.650	ml/mol	McGowan Method
rinpol	1501.00		NIST Webbook

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R119056&Units=SI

Legend

log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
rinpol:	Non-polar retention indices

Latest version available from:

<https://www.chemeo.com/cid/37-768-9/Methanimine-1-1-pyrrolidiny-N-hexyl.pdf>

Generated by Cheméo on 2025-12-22 15:54:10.402041 +0000 UTC m=+6167047.932081653.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.